

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/13/2002 Time: 07:49:38

Sample: AD31363
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/13/02 07:49 Ended: 2/13/02 07:49 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		95		5.0

Comments for Sample AD31363 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-13-2002 Time: 07:50:26

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

The recoveries for 3 of the 43 compounds in the LCS Standard were above their acceptance ranges.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\JAN1602\31363.D

Vial: 8

Acq On : 16 Jan 2002 6:29 pm

Operator: 209 GALOS

Sample : gc/ms scan 12/27

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Feb 7 8:43 2002

Quant Results File: GCMS0103.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Mon Jan 07 10:31:02 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0103

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.50	152	804402	20.00	ug/l	-0.03
10) Naphthalene-d8	15.09	136	3140872	20.00	ug/l	-0.02
17) Acenaphthene-d10	20.35	164	1556008	20.00	ug/l	-0.03
29) Phenanthrene-d10	24.77	188	2038965	20.00	ug/l	-0.03
38) Chrysene-d12	32.80	240	1126838	20.00	ug/l	-0.03
44) Perylene-d12	36.86	264	698976	20.00	ug/l	-0.03

System Monitoring Compounds

37) 2,4-DDD	29.85	235	110588	4.73	ug/mL	-0.22
Spiked Amount	5.000		Recovery	=	94.60%	

Target Compounds

					Qvalue
2) N-nitroso-dimethyl amine	0.00	74	0	N.D.	
3) bis(2-Chloroethyl)ether	0.00	93	0	N.D.	
4) 1,3-Dichlorobenzene	0.00	146	0	N.D.	
5) 1,4-Dichlorobenzene	0.00	146	0	N.D.	
6) 1,2-Dichlorobenzene	0.00	146	0	N.D.	
7) 2,2'-oxybis(1-Chloropropan	0.00	45	0	N.D.	
8) N-Nitroso-di-n-propylamine	0.00	70	0	N.D.	
9) Hexachloroethane	0.00	117	0	N.D.	
11) Nitrobenzene	0.00	77	0	N.D.	
12) Isophorone	13.32	82	297	N.D.	
13) bis(2-Chloroethoxy)methane	0.00	93	0	N.D.	
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.	
15) Naphthalene	15.15	128	884	N.D.	
16) Hexachlorobutadiene	0.00	225	0	N.D.	
18) Hexachlorocyclopentadiene	0.00	237	0	N.D.	
19) 2-Chloronaphthalene	0.00	162	0	N.D.	
20) Dimethylphthalate	0.00	163	0	N.D.	
21) 2,6-Dinitrotoluene	0.00	165	0	N.D.	d
22) Acenaphthylene	0.00	152	0	N.D.	
23) Acenaphthene	0.00	153	0	N.D.	
24) 2,4-Dinitrotoluene	20.84	165	187	N.D.	
25) Diethylphthalate	21.89	149	287	N.D.	
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.	
27) Fluorene	0.00	166	0	N.D.	
28) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.	

(#)=qualifier out of range (m)=manual integration

31363.D GCMS0103.M

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Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\JAN1602\31363.D

Vial: 8

Acq On : 16 Jan 2002 6:29 pm

Operator: 209 GALOS

Sample : gc/ms scan 12/27

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Feb 7 8:43 2002

Quant Results File: GCMS0103.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Mon Jan 07 10:31:02 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0103

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
30) N-Nitrosodiphenylamine	0.00	168	0	N.D.		
31) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		
32) Hexachlorobenzene	0.00	284	0	N.D.		
33) Phenanthrene	24.77	178	704	N.D.		
34) Anthracene	24.77	178	704	N.D.		
35) Di-n-butylphthalate	26.80	149	26048	N.D.		
36) Fluoranthene	0.00	202	0	N.D.		
39) Pyrene	0.00	202	0	N.D.		
40) Butylbenzylphthalate	0.00	149	0	N.D.		
41) Benzo(a)anthracene	32.80	228	2839	N.D.		
42) Bis(2-ethylhexyl)phthalate	33.10	149	1270	N.D.		
43) Chrysene	32.80	228	2839	N.D.		
45) Di-n-Octylphthalate	0.00	149	0	N.D.		
46) Benzo(b)fluoranthene	0.00	252	0	N.D.		
47) Benzo(k)fluoranthene	0.00	252	0	N.D.		
48) Benzo(a)pyrene	36.86	252	2712	N.D.		
49) Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.		
50) Dibenz(a,h)anthracene	0.00	278	0	N.D.		
51) Benzo(g,h,i)perylene	0.00	276	0	N.D.		

(#) = qualifier out of range (m) = manual integration

31363.D GCMS0103.M

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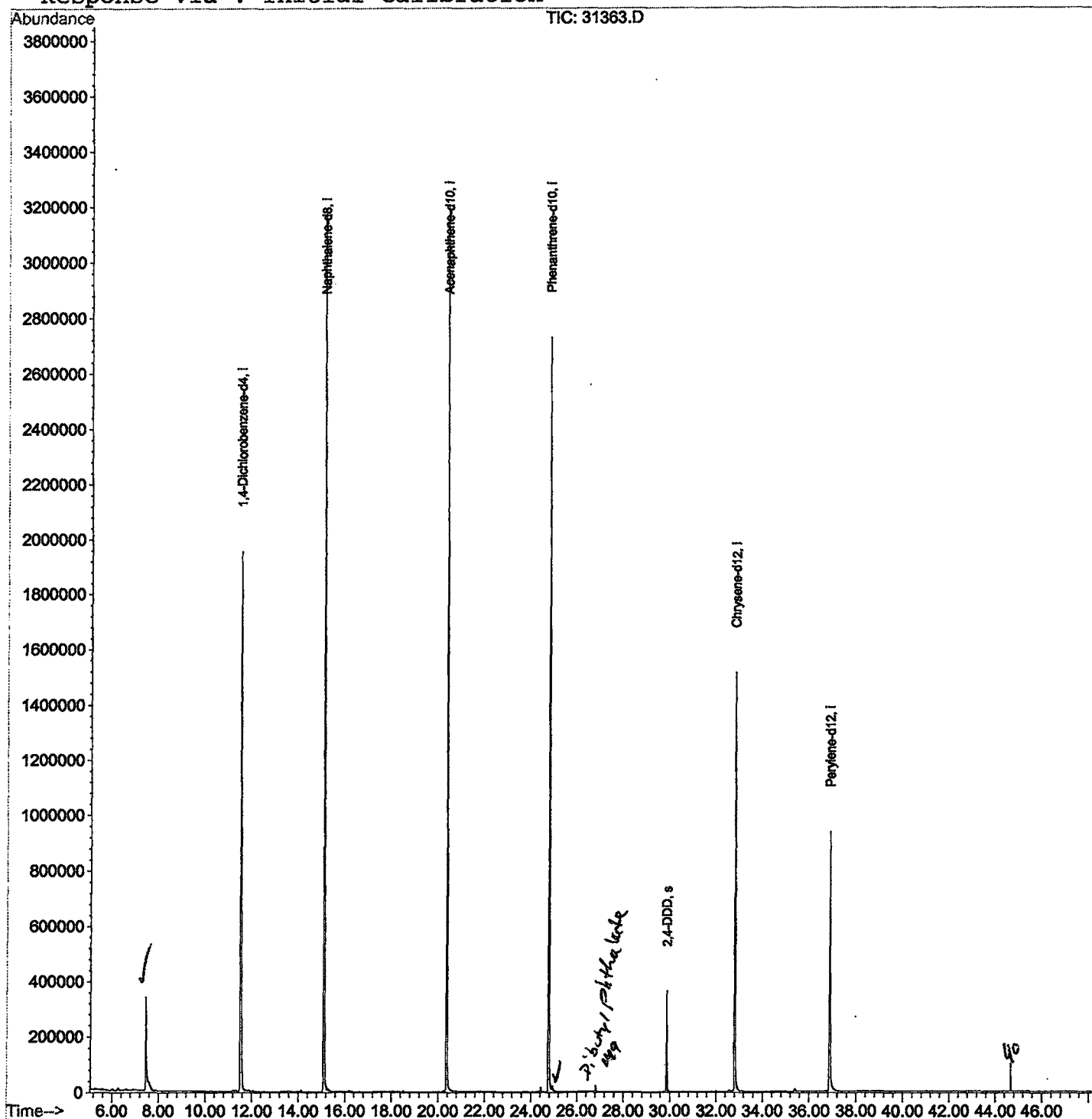
Quantitation Report

Data File : C:\HPCHEM\1\DATA\JAN1602\31363.D
Acq On : 16 Jan 2002 6:29 pm
Sample : gc/ms scan 12/27
Misc :
MS Integration Params: GOOFY.P
Quant Time: Feb 7 8:43 2002

Vial: 8
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Results File: GCMS0103.RES

Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Mon Jan 07 10:31:02 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\JAN1602\31363.D

Acq On : 16 Jan 2002 6:29 pm

Sample : gc/ms scan 12/27

Misc :

MS Integration Params: LSCINT.P

Vial: 8

Operator: 209 GALOS

Inst : GC/MS 209

Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.447	258	262	275	rBV	335768	884261	13.41%	2.778%
2	11.496	699	703	723	rBV	1952303	5101388	77.37%	16.026%
3	15.095	1089	1095	1114	rBV	3174048	6463214	98.02%	20.304%
4	20.355	1662	1668	1688	rBV	3205019	6593692	100.00%	20.714%
5	24.771	2143	2149	2162	rBV2	2729239	5780222	87.66%	18.158%
6	24.918	2162	2165	2182	rVB	21199	70232	1.07%	0.221%
7	29.847	2697	2702	2712	rVB2	361900	703149	10.66%	2.209%
8	32.803	3017	3024	3043	rBV2	1511589	3640444	55.21%	11.436%
9	36.861	3459	3466	3486	rBV	933334	2595956	39.37%	8.155%

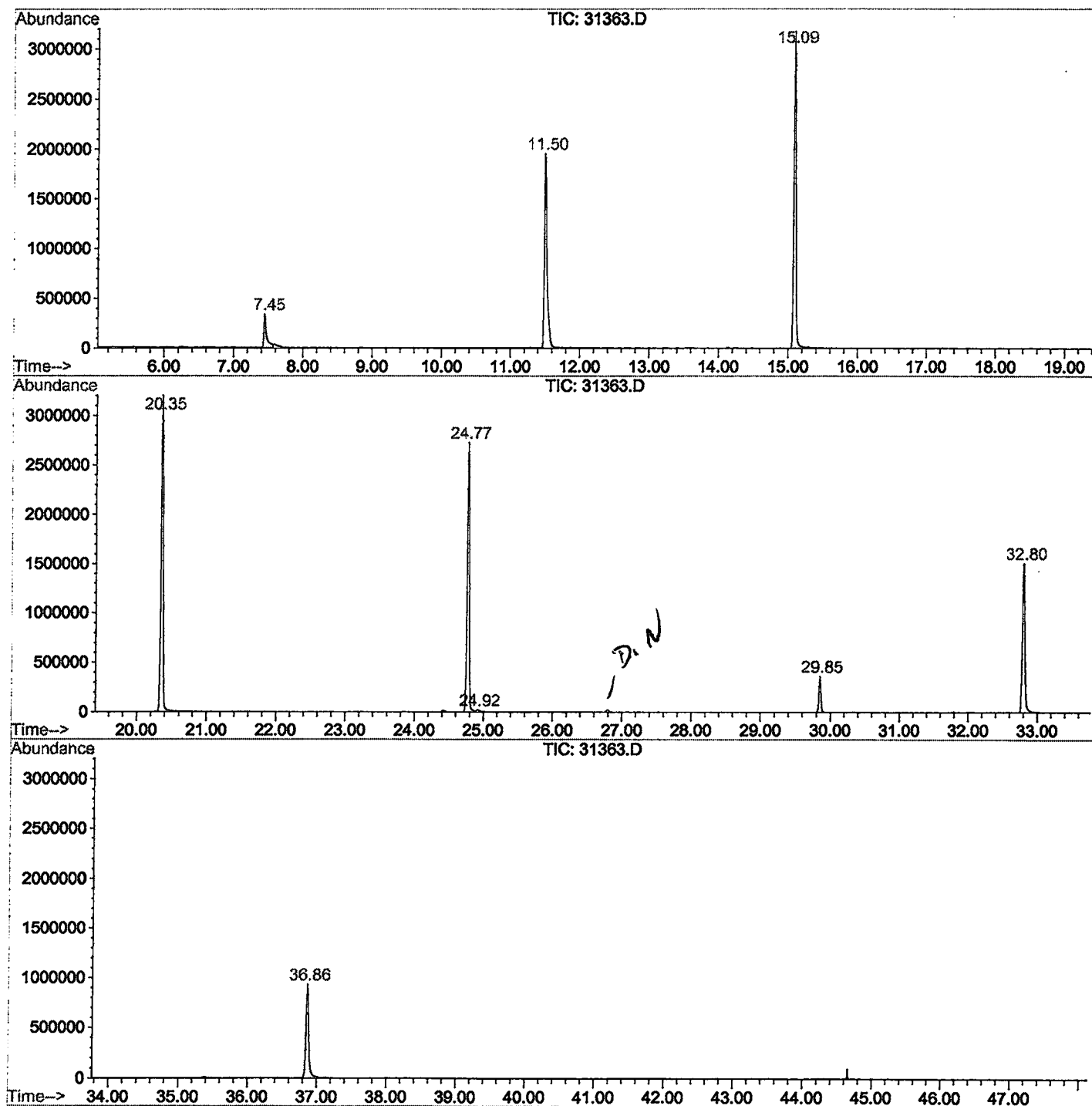
Sum of corrected areas: 31832558

31363.D GCMS0103.M

Thu Feb 07 09:11:11 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\JAN1602\31363.D
Operator : 209 GALOS
Acquired : 16 Jan 2002 6:29 pm using AcqMethod GCMS0103
Instrument : GC/MS 209
Sample Name: gc/ms scan 12/27
Misc Info :
Vial Number: 8
Quant File :GCMS0103.RES (RTE Integrator)



31363.D GCMS0103.M

Thu Feb 07 09:11:17 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN1602\31363.D
 Acq On : 16 Jan 2002 6:29 pm
 Sample : gc/ms scan 12/27
 Misc :
 MS Integration Params: LSCINT.P

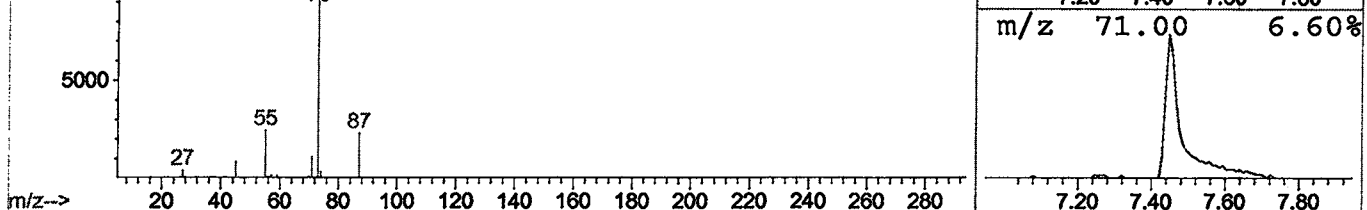
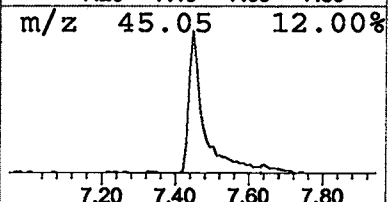
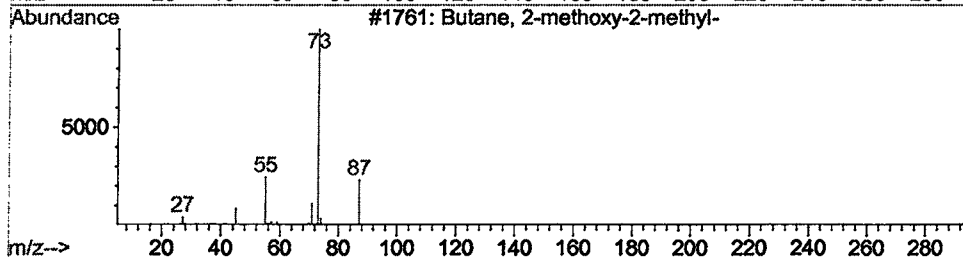
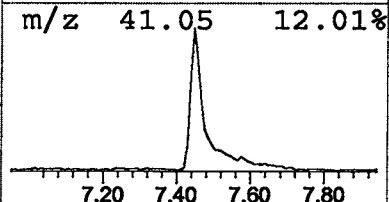
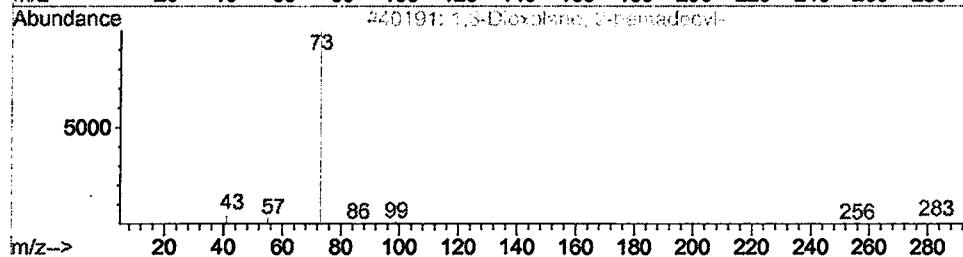
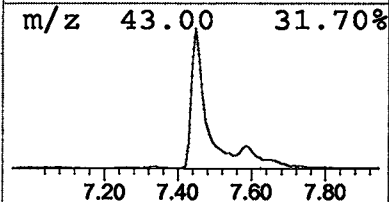
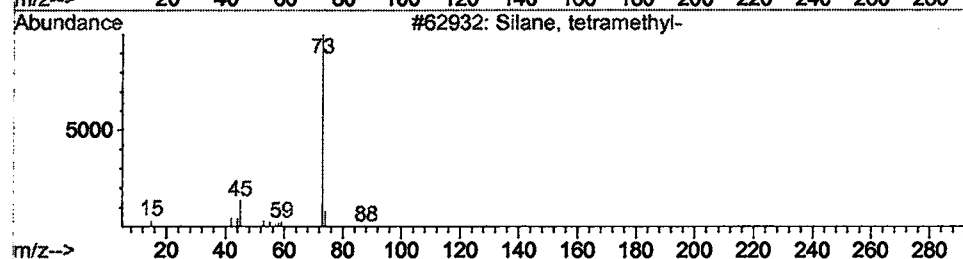
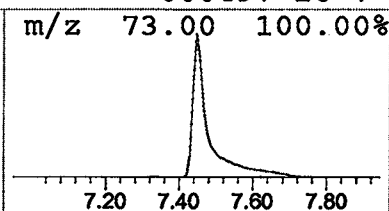
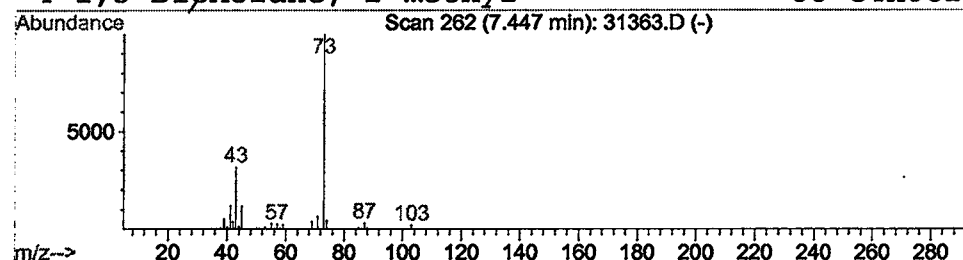
Vial: 8
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Library : C:\DATABASE\NBS75K.L

 Peak Number 1 Silane, tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.45	3.47 ug/l	884261	1,4-Dichlorobenzene-d4	11.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9	
2	1,3-Dioxolane, 2-pentadecyl-	284	C18H36O2	004360-57-0	9	
3	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9	
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	7	



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN1602\31363.D
 Acq On : 16 Jan 2002 6:29 pm
 Sample : gc/ms scan 12/27
 Misc :
 MS Integration Params: LSCINT.P

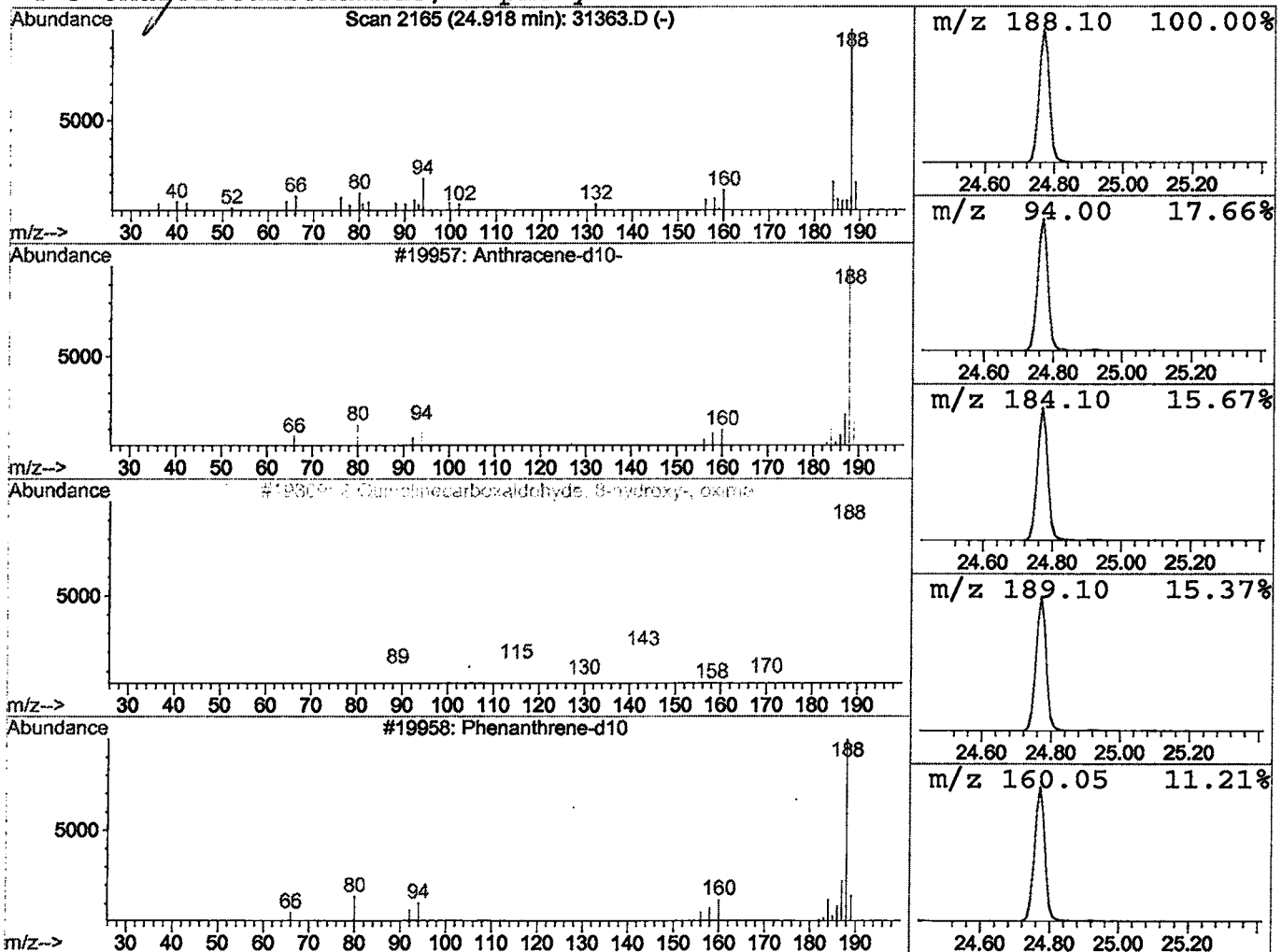
Vial: 8
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Library : C:\DATABASE\NBS75K.L

 Peak Number 2 Anthracene-d10- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.92	0.24 ug/l	70232	Phenanthrene-d10	24.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene-d10- 5	188	C14D10	001719-06-8	91
2			2-Quinolincarboxaldehyde, 8-hydrox	188	C10H8N2O2	005603-22-5	38
3			Phenanthrene-d10	188	C14D10	001517-22-2	38
4			5-Oxazolecarboxamide, 2-phenyl-	188	C10H8N2O2	039819-42-6	38



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 16 Jan 2002 6:29 pm
Data File: C:\HPCHEM\1\DATA\JAN1602\31363.D
Name: gc/ms scan 12/27
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title: 209 625/TCLP calibration table
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Silane, tetramethyl-	7.45	3.5	ug/l	884261	ISTD01	11.50	5101390	20.0
Anthracene-d10- SS	24.92	0.2	ug/l	70232	ISTD04	24.77	5780220	20.0

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